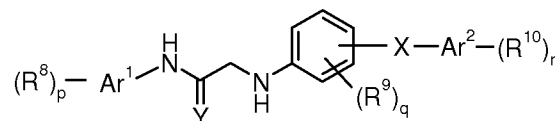


This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (Cancelled):
2. (Cancelled):
3. (Currently Amended): A glycine compound according to ~~claim 1~~, selected from the compounds of formula II: [[,]]



II

wherein

Ar¹ is phenyl, pyridinyl, oxazolyl, isoxazolyl, pyrazolyl or imidazolyl;

Ar¹, Ar² are each, independently from one another, selected from is an aromatic hydrocarbon group hydrocarbons containing 6 to 14 carbon atoms or an and ethylenical unsaturated or aromatic heterocyclic group residues containing 3 to 10 carbon atoms and one or two hetero atoms, wherein said hetero atoms are each independently selected from N, O or and S; [[,]]

R⁸ is alkyl having 1 to 4 carbon atoms, alkoxy having 1 to 4 carbon atoms, Hal, CH₂Hal, CH(Hal)₂, perhaloalkyl having 1 to 4 carbon atoms, NO₂, (CH₂)_nCN, (CH₂)_nNR¹¹R¹², (CH₂)_nCOR¹³, (CH₂)_nCOOR¹¹, (CH₂)_nCONR¹¹R¹², (CH₂)_nSO₂NR¹¹R¹², or (CH₂)_nS(O)_uR¹³;

R⁸, R⁹ and R¹⁰ are each independently selected from H, A, cycloalkyl having

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~~comprising~~ 3 to 7 carbon atoms, Hal, CH₂Hal, CH(Hal)₂, C(Hal)₃, NO₂,
 (CH₂)_nCN, (CH₂)_nNR¹¹R¹², (CH₂)_nOR¹¹, (CH₂)_nO(CH₂)_kNR¹¹R¹²,
 (CH₂)_nCOOR¹², (CH₂)_nCONR¹¹R¹², (CH₂)_nNR¹¹COR¹³,
 (CH₂)_nNR¹¹CONR¹¹R¹², (CH₂)_nNR¹¹SO₂A, (CH₂)_nSO₂NR¹¹R¹²,
 (CH₂)_nS(O)_uR¹³, (CH₂)_nOC(O)R¹³, (CH₂)_nCOR¹³, (CH₂)_nSR¹¹, CH=N-OA,
 CH₂CH=N-OA, (CH₂)_nNHOA, (CH₂)_nCH=N-R¹¹, (CH₂)_nOC(O)NR¹¹R¹²,
 (CH₂)_nNR¹¹COOR¹², (CH₂)_nN(R¹¹)CH₂CH₂OR¹³, (CH₂)_nN(R¹¹)CH₂CH₂OCF₃,
 (CH₂)_nN(R¹¹)C(R¹³)HCOOR¹², C(R¹³)HCOOR¹²,
 (CH₂)_nN(R¹¹)CH₂CH₂N(R¹²)CH₂COOR¹², (CH₂)_nN(R¹¹)CH₂CH₂NR¹¹R¹²,
 CH=CHCOOR¹¹, CH=CHCH₂NR¹¹R¹², CH=CHCH₂NR¹¹R¹²,
 CH=CHCH₂OR¹³, (CH₂)_nN(COOR¹¹)COOR¹², (CH₂)_nN(CONH₂)COOR¹¹,
 (CH₂)_nN(CONH₂)CONH₂, (CH₂)_nN(CH₂COOR¹¹)COOR¹²,
 (CH₂)_nN(CH₂CONH₂)COOR¹¹, (CH₂)_nN(CH₂CONH₂)CONH₂,
 (CH₂)_nCHR¹³COR¹¹, (CH₂)_nCHR¹³COOR¹¹, (CH₂)_nCHR¹³CH₂OR¹⁴,
 (CH₂)_nOCN₂or ~~and~~ (CH₂)_nNCO₂; [[,]]

R¹¹, R¹² are each independently ~~selected from~~ H, A, (CH₂)_mAr³, or ~~and~~ (CH₂)_mHet, or,
 in NR¹¹R¹², R¹¹ and R¹² form, together with the N-Atom they are bound to, a 5-, 6- or
 7-membered heterocycle which optionally contains 1 or 2 additional hetero
 atoms, wherein said hetero atoms are each independently selected from N, O
or an S_i; [[,]]

R¹³, R¹⁴ are each independently ~~selected from~~ H, Hal, A, (CH₂)_mAr⁴, or ~~and~~
 (CH₂)_mHet_i; [[,]]

A is ~~selected from~~ alkyl, alkenyl, cycloalkyl, alkylencycloalkyl, alkoxy or ~~and~~
 alkoxyalkyl_i; [[,]]

Ar³, Ar⁴ are each independently from one another an aromatic hydrocarbon group
having residues comprising 5 to 12 which is ~~are~~ optionally substituted by one
or more substituents, wherein said substituents are in each case independently

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~~selected from~~ A, Hal, NO₂, CN, OR¹⁵, NR¹⁵R¹⁶, COOR¹⁵, CONR¹⁵R¹⁶,
NR¹⁵COR¹⁶, NR¹⁵CONR¹⁵R¹⁶, NR¹⁶SO₂A, COR¹⁵, SO₂R¹⁵R¹⁶, S(O)_uA, or
~~and~~ OOCR¹⁵; [[,]]

Het is a saturated, unsaturated or aromatic heterocyclic group residue which is
optionally substituted by one or more substituents, wherein said substituents
are in each case independently selected from A, Hal, NO₂, CN, OR¹⁵, NR¹⁵R¹⁶,
COOR¹⁵, CONR¹⁵R¹⁶, NR¹⁵COR¹⁶, NR¹⁵CONR¹⁵R¹⁶, NR¹⁶SO₂A, COR¹⁵,
SO₂R¹⁵R¹⁶, S(O)_uA, or ~~and~~ OOCR¹⁵; [[,]]

R¹⁵, R¹⁶ are each independently ~~selected from~~ H, A, or ~~and~~ (CH₂)_mAr⁵; [[,]]

Ar⁵ Ar⁶ is a 5- or 6-membered aromatic hydrocarbon which is optionally substituted
by one or more substituents, wherein said substituents are in each case
independently selected from methyl, ethyl, propyl, 2-propyl, tert.-butyl, Hal,
CN, OH, NH₂, or ~~and~~ CF₃; [[,]]

k, ~~n~~, m are independently of one another 0, 1, 2, 3, 4, or 5;

n is 0 or 1;

X is O, S, NR¹⁵, CHOR¹¹, CH₂, CH₂CH₂, OCH₂, CH₂O, OCH₂CH₂, or
CH₂CH₂O; represents a bond or is (CR¹¹R¹²)_h, or (CHR¹¹)_h-Q-(CHR¹²)_i;

Q ~~is selected from~~ O, S, N-R¹⁵, (CHAl₂)_j, (O-CHR¹⁸)_j, (CHR¹⁸-O)_j, CR¹⁸=CR¹⁹,
(O-CHR¹⁸CHR¹⁹)_j, CHR¹⁸CHR¹⁹-O)_j, C=O, C=S, C=NR¹⁵, CH(OR¹⁵),
C(OR¹⁷)(OR²⁰), C(=O)O, OC(=O), OC(=O)O, C(=)N(R¹⁵), N(R¹⁵)C(=O),
OC(=O)N(R¹⁵), N(R¹⁵)C(=O)O, CH=N-O, CH=N-NR¹⁵, OC(O)NR¹⁵,
NR¹⁵C(O)O, S=O, SO₂, SO₂NR¹⁵, or ~~and~~ NR¹⁵SO₂; [[,]]

h, i are independently from each other 0, 1, 2, 3, 4, 5 or 6; [[,]]

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j is 1, 2, 3, 4, 5 or 6; [[,]]

Y is ~~selected from~~ O, S, NR²¹, C(R²²)-NO₂, C(R²²)-CN, or ~~and~~ C(CN)₂; [[,]]

R²¹ is H, Hal, A, (CH₂)_mAr⁴, or ~~and~~ (CH₂)_mHet_i; [[,]]

R²² is H, A, (CH₂)_mAr³, or ~~and~~ (CH₂)_mHet_i; [[,]]

p is 1, 2 or 3;

~~p, r~~ is ~~are independently from one another~~ 0, 1, 2, 3, 4 or 5; [[,]]

q is 0 or 1; ~~0, 1, 2, 3 or 4~~,

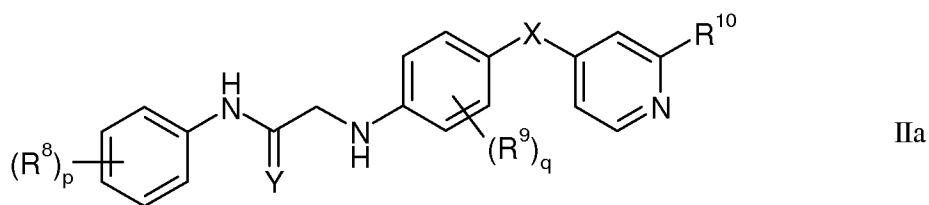
u is 0 or 2; ~~0, 1, 2 or 3~~,

and

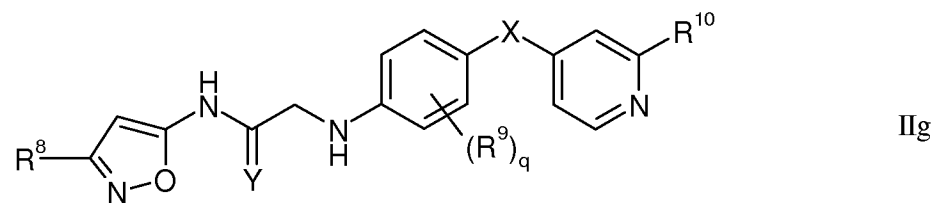
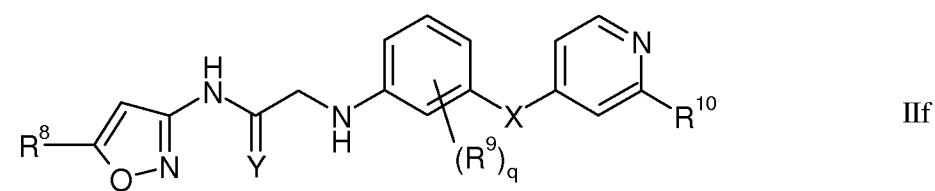
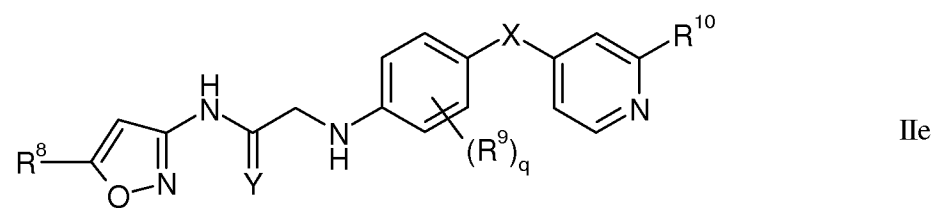
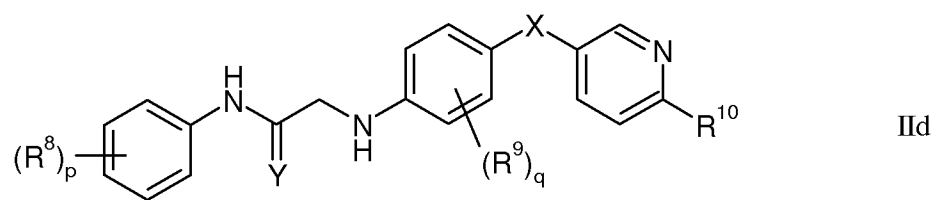
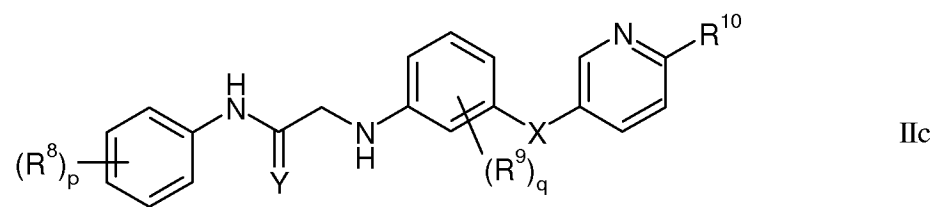
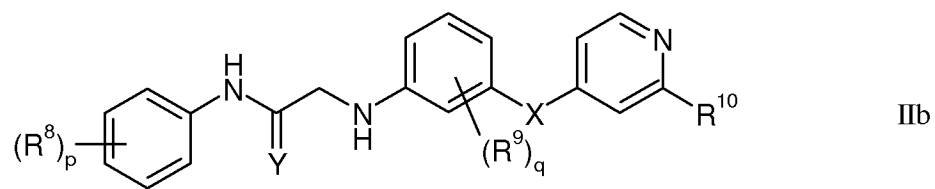
Hal is in each case independently ~~selected from~~ F, Cl, Br or ~~and~~ I;

or a salt or solvate ~~and the salts and solvates~~ thereof.

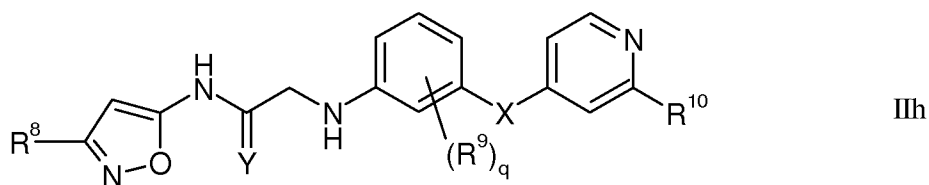
4. (Currently Amended): A glycine compound according to claim 3 ~~4~~,
selected from the compounds of formula IIa, IIb, IIc, IId, IIe, IIg and IIh,
[[']]



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wherein

R⁸ is alkyl having 1 to 4 carbon atoms, alkoxy having 1 to 4 carbon atoms, Hal, CH₂Hal, CH(Hal)₂, C(Hal)₃, NO₂, (CH₂)_nCN, (CH₂)_nNR¹¹R¹², (CH₂)_nCOR¹³, (CH₂)_nCOOR¹¹, (CH₂)_nCONR¹¹R¹², (CH₂)_nSO₂NR¹¹R¹², or (CH₂)_nS(O)_uR¹³,

R⁸, R⁹ and R¹⁰ are independently selected from H, A, cycloalkyl comprising 3 to 7 carbon atoms, Hal, CH₂Hal, CH(Hal)₂, C(Hal)₃, NO₂, (CH₂)_nCN, (CH₂)_nNR¹¹R¹², (CH₂)_nOR¹¹, (CH₂)_nO(CH₂)_kNR¹¹R¹², (CH₂)_nCOOR¹², (CH₂)_nCONR¹¹R¹², (CH₂)_nNR¹¹COR¹³, (CH₂)_nNR¹¹CONR¹¹R¹², (CH₂)_nNR¹¹SO₂A, (CH₂)_nSO₂NR¹¹R¹², (CH₂)_nS(O)_uR¹³, (CH₂)_nOC(O)R¹³, (CH₂)_nCOR¹³, (CH₂)_nSR¹¹, CH=N-OA, CH₂CH=N-OA, (CH₂)_nNHOA, (CH₂)_nCH=N-R¹¹, (CH₂)_nOC(O)NR¹¹R¹², (CH₂)_nNR¹¹COOR¹², (CH₂)_nN(R¹¹)CH₂CH₂OR¹³, (CH₂)_nN(R¹¹)CH₂CH₂OCF₃, (CH₂)_nN(R¹¹)C(R¹³)HCOOR¹², C(R¹³)HCOOR¹², (CH₂)_nN(R¹¹)CH₂CH₂N(R¹²)CH₂COOR¹², (CH₂)_nN(R¹¹)CH₂CH₂NR¹¹R¹², CH=CHCOOR¹¹, CH=CHCH₂NR¹¹R¹², CH=CHCH₂NR¹¹R¹², CH=CHCH₂OR¹³, (CH₂)_nN(COOR¹¹)COOR¹², (CH₂)_nN(CONH₂)COOR¹¹, (CH₂)_nN(CONH₂)CONH₂, (CH₂)_nN(CH₂COOR¹¹)COOR¹², (CH₂)_nN(CH₂CONH₂)COOR¹¹, (CH₂)_nN(CH₂CONH₂)CONH₂, (CH₂)_nCHR¹³COR¹¹, (CH₂)_nCHR¹³COOR¹¹, (CH₂)_nCHR¹³CH₂OR¹⁴, (CH₂)_nOCN and (CH₂)_nNCO,

R¹⁰ can also be H,

R¹¹, R¹² are each independently ~~selected from~~ H, A, (CH₂)_mAr³, or and (CH₂)_mHet, or, in NR¹¹R¹², R¹¹ and R¹² form, together with the N-Atom they are bound to, a 5-, 6- or 7-membered heterocycle which optionally contains 1 or 2 additional

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hetero atoms, wherein said hetero atoms are independently selected from N, O or an S,

R^{13} , R^{14} are each independently ~~selected from~~ H, Hal, A, $(CH_2)_m Ar^4$, or and $(CH_2)_m Het$,

A is ~~selected from~~ alkyl, alkenyl, cycloalkyl, alkylencycloalkyl, alkoxy, or and alkoxyalkyl,

Ar^3 , Ar^4 are each independently from one another an aromatic hydrocarbon group having residues comprising 5 to 12 carbon atoms which is are optionally substituted by one or more substituents, wherein said substituents are in each case independently selected from A, Hal, NO_2 , CN, OR^{15} , $NR^{15}R^{16}$, $COOR^{15}$, $CONR^{15}R^{16}$, $NR^{15}COR^{16}$, $NR^{15}CONR^{15}R^{16}$, $NR^{16}SO_2A$, COR^{15} , $SO_2R^{15}R^{16}$, $S(O)_uA$, or and $OOCR^{15}$,

Het is a saturated, unsaturated or aromatic heterocyclic group residue which is optionally substituted by one or more substituents, wherein said substituents are in each case independently selected from A, Hal, NO_2 , CN, OR^{15} , $NR^{15}R^{16}$, $COOR^{15}$, $CONR^{15}R^{16}$, $NR^{15}COR^{16}$, $NR^{15}CONR^{15}R^{16}$, $NR^{16}SO_2A$, COR^{15} , $SO_2R^{15}R^{16}$, $S(O)_uA$, or and $OOCR^{15}$,

R^{15} , R^{16} are each independently ~~selected from a group consisting of~~ H, A, or and $(CH_2)_m Ar^5$,

Ar^5 is a 5- or 6-membered aromatic hydrocarbon which is optionally substituted by one or more substituents, wherein said substituents are in each case independently selected from methyl, ethyl, propyl, 2-propyl, tert.-butyl, Hal, CN, OH, NH_2 , or and CF_3 ,

k, ~~n~~, m are independently of one another 0, 1, 2, 3, 4, or 5;

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p is 0, 1, or 2, ~~0, 1, 2, 3, 4 or 5~~,

q is 0 or 1, ~~0, 1, 2, 3 or 4~~,

u is 0, 1, 2 or 3,

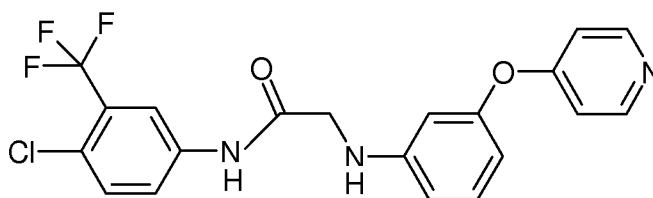
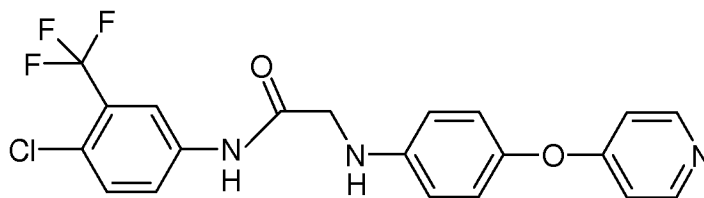
Y is ~~selected from~~ O, S, NR²¹, C(R²²)-NO₂, C(R²²)-CN, or ~~and~~ C(CN)₂,

R²¹ is H, Hal, A, (CH₂)_mAr⁴, or ~~and~~ (CH₂)_mHet,

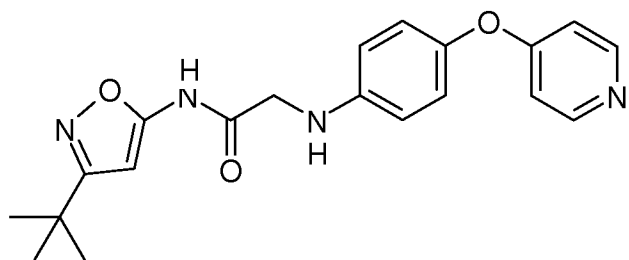
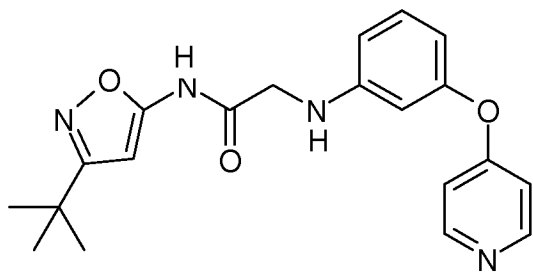
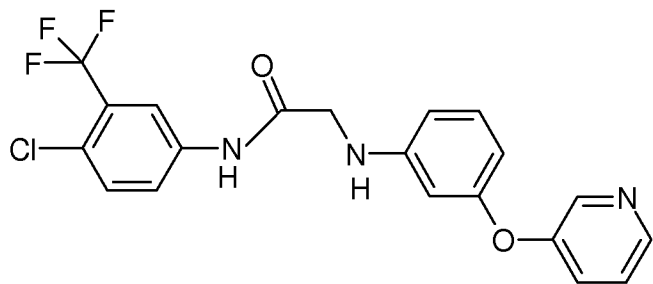
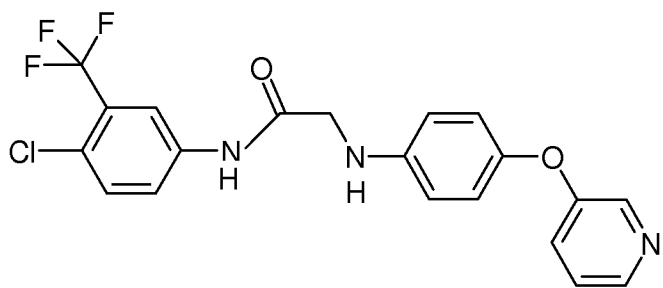
R²² is H, A, (CH₂)_mAr³, or ~~and~~ (CH₂)_mHet,

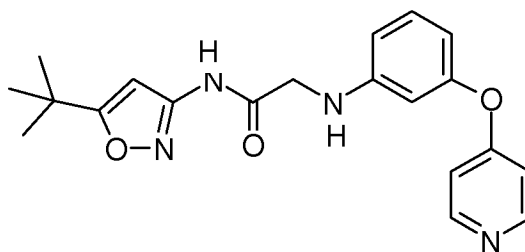
or a salt or solvate ~~and the salts and solvates~~ thereof.

5. (Currently Amended): A glycineamide compound according to claim 4 3,
selected from

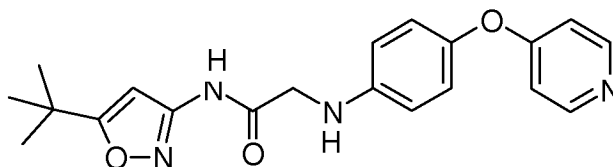


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and



6. (Cancelled):

7. (Cancelled):

8. ((Cancelled):

9. (Cancelled):

10. (Previously Presented): A pharmaceutical composition comprising one or more compounds according to claim 1, and one or more additional compounds, selected from physiologically acceptable excipients, auxiliaries, adjuvants, carriers and other pharmaceutical active ingredients.

11. (Currently Amended): A process for the manufacture of a pharmaceutical composition comprising:

processing one or more compounds according to claim 1 and one or more compounds selected from the group consisting of carriers, excipients, auxiliaries and pharmaceutical active ingredients other than the compounds according to claim 1, by mechanical means into a dosage form pharmaceutical composition that is suitable as ~~dosage form~~ for application

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and/or administration to a patient.

12. (Cancelled):

13. (Cancelled):

14. (Cancelled):

15. (Cancelled):

16. (Cancelled):

17. (Cancelled):

18. (Cancelled):

19. (Cancelled):

20. (Cancelled):

21. (Cancelled):

22. (Cancelled):

23. (Cancelled):

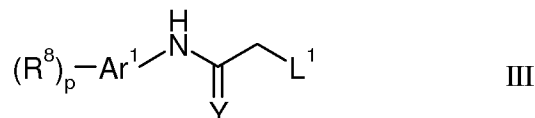
24. (Cancelled):

25. (Cancelled):

26. (Cancelled):

27. (Cancelled):

28. (Currently Amended): A method for producing a compound according to claim 3 of ~~formula II~~, said method comprising:
reacting a compound of formula III



wherein

L^1 is Cl, Br, I, OH, a reactive esterified OH-group or a diazonium moiety, and

R^8 , Ar^1 , and Y are as defined in claim 3,

R^8 is selected from H, A, cycloalkyl comprising 3 to 7 carbon atoms, Hal, CH_2Hal , $\text{CH}(\text{Hal})_2$, $\text{C}(\text{Hal})_3$, NO_2 , $(\text{CH}_2)_n\text{CN}$, $(\text{CH}_2)_n\text{NR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{OR}^{11}$, $(\text{CH}_2)_n\text{O}(\text{CH}_2)_k\text{NR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{COOR}^{12}$, $(\text{CH}_2)_n\text{CONR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{NR}^{11}\text{COR}^{13}$, $(\text{CH}_2)_n\text{NR}^{11}\text{CONR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{NR}^{11}\text{SO}_2\text{A}$, $(\text{CH}_2)_n\text{SO}_2\text{NR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{S}(\text{O})_u\text{R}^{13}$, $(\text{CH}_2)_n\text{OC}(\text{O})\text{R}^{13}$, $(\text{CH}_2)_n\text{COR}^{13}$, $(\text{CH}_2)_n\text{SR}^{11}$, $\text{CH}=\text{N}-\text{OA}$, $\text{CH}_2\text{CH}=\text{N}-\text{OA}$, $(\text{CH}_2)_n\text{NHOA}$, $(\text{CH}_2)_n\text{CH}=\text{N}-\text{R}^{11}$, $(\text{CH}_2)_n\text{OC}(\text{O})\text{NR}^{11}\text{R}^{12}$, $(\text{CH}_2)_n\text{NR}^{11}\text{COOR}^{12}$, $(\text{CH}_2)_n\text{N}(\text{R}^{11})\text{CH}_2\text{CH}_2\text{OR}^{13}$, $(\text{CH}_2)_n\text{N}(\text{R}^{11})\text{CH}_2\text{CH}_2\text{OCF}_3$, $(\text{CH}_2)_n\text{N}(\text{R}^{11})\text{C}(\text{R}^{13})\text{HCOOR}^{12}$, $\text{C}(\text{R}^{13})\text{HCOOR}^{12}$, $(\text{CH}_2)_n\text{N}(\text{R}^{11})\text{CH}_2\text{CH}_2\text{N}(\text{R}^{12})\text{CH}_2\text{COOR}^{12}$, $(\text{CH}_2)_n\text{N}(\text{R}^{11})\text{CH}_2\text{CH}_2\text{NR}^{11}\text{R}^{12}$, $\text{CH}=\text{CHCOOR}^{11}$, $\text{CH}=\text{CHCH}_2\text{NR}^{11}\text{R}^{12}$, $\text{CH}=\text{CHCH}_2\text{NR}^{11}\text{R}^{12}$, $\text{CH}=\text{CHCH}_2\text{OR}^{13}$, $(\text{CH}_2)_n\text{N}(\text{COOR}^{11})\text{COOR}^{12}$, $(\text{CH}_2)_n\text{N}(\text{CONH}_2)\text{COOR}^{11}$, $(\text{CH}_2)_n\text{N}(\text{CONH}_2)\text{CONH}_2$, $(\text{CH}_2)_n\text{N}(\text{CH}_2\text{COOR}^{11})\text{COOR}^{12}$, $(\text{CH}_2)_n\text{N}(\text{CH}_2\text{CONH}_2)\text{COOR}^{11}$, $(\text{CH}_2)_n\text{N}(\text{CH}_2\text{CONH}_2)\text{CONH}_2$,

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$(\text{CH}_2)_n\text{CHR}^{13}\text{COR}^{11}$, $(\text{CH}_2)_n\text{CHR}^{13}\text{COOR}^{14}$, $(\text{CH}_2)_n\text{CHR}^{13}\text{CH}_2\text{OR}^{14}$,
 $(\text{CH}_2)_n\text{OCN}$ and $(\text{CH}_2)_n\text{NCO}$,

p — is 0, 1, 2, 3, 4 or 5,

Ar^1 — is selected from aromatic hydrocarbons containing 6 to 14 carbon atoms and ethylenical unsaturated or aromatic heterocyclic residues containing 3 to 10 carbon atoms and one or two hetero atoms, independently selected from N, O and S,

Y — is selected from O, S, NR^{21} , $\text{C}(\text{R}^{22})\text{NO}_2$, $\text{C}(\text{R}^{22})\text{CN}$ and $\text{C}(\text{CN})_2$,

A — is selected from alkyl, alkenyl, cycloalkyl, alkylencycloalkyl, alkoxy and alkoxyalkyl,

R^{11} , R^{12} are independently selected from H, A, $(\text{CH}_2)_m\text{Ar}^3$ and $(\text{CH}_2)_m\text{Het}$, or, in $\text{NR}^{11}\text{R}^{12}$, R^{11} and R^{12} form, together with the N Atom they are bound to, a 5, 6 or 7 membered heterocycle which optionally contains 1 or 2 additional hetero atoms, selected from N, O and S,

k, n, m are independently of one another 0, 1, 2, 3, 4, or 5,

R^{13} , R^{14} are independently selected from H, Hal, A, $(\text{CH}_2)_m\text{Ar}^4$ and $(\text{CH}_2)_m\text{Het}$,

Ar^3 , Ar^4 are independently from one another aromatic hydrocarbon residues comprising 5 to 12 carbon atoms which are optionally substituted by one or more substituents, selected from A, Hal, NO_2 , CN, OR^{15} , $\text{NR}^{15}\text{R}^{16}$, COOR^{15} , $\text{CONR}^{15}\text{R}^{16}$, $\text{NR}^{15}\text{COR}^{16}$, $\text{NR}^{15}\text{CONR}^{15}\text{R}^{16}$, $\text{NR}^{16}\text{SO}_2\text{A}$, COR^{15} , $\text{SO}_2\text{R}^{15}\text{R}^{16}$, $\text{S}(\text{O})_u\text{A}$ and OOCR^{15} ,

Het — is a saturated, unsaturated or aromatic heterocyclic residue which is

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~~optionally substituted by one or more substituents, selected from A, Hal, NO₂, CN, OR¹⁵, NR¹⁵R¹⁶, COOR¹⁵, CONR¹⁵R¹⁶, NR¹⁵COR¹⁶, NR¹⁵CONR¹⁵R¹⁶, NR¹⁶SO₂A, COR¹⁵, SO₂R¹⁵R¹⁶, S(O)_uA and OOCR¹⁵;~~

~~R¹⁵, R¹⁶ are independently selected from H, A, and (CH₂)_mAr⁵;~~

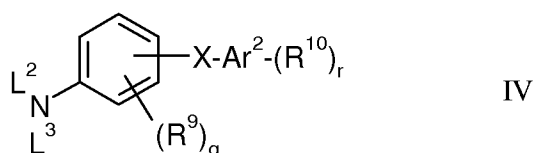
~~Ar⁵ is a 5 or 6 membered aromatic hydrocarbon which is optionally substituted by one or more substituents selected from methyl, ethyl, propyl, 2-propyl, tert. butyl, Hal, CN, OH, NH₂ and CF₃;~~

~~u is 0, 1, 2 or 3;~~

~~R²⁴ is H, Hal, A, (CH₂)_mAr⁴ and (CH₂)_mHet;~~

~~R²² is H, A, (CH₂)_mAr³ and (CH₂)_mHet;~~

with a compound of formula IV,



wherein

L², L³ are independently from one another H or a metal ion, and

R⁹, q, X, Ar², R¹⁰ and r are as defined in claim 3,

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R^9 and R^{10} are independent from one another selected from H, A, cycloalkyl comprising 3 to 7 carbon atoms, Hal, CH_2Hal , $CH(Hal)_2$, $C(Hal)_3$, NO_2 , $(CH_2)_nCN$, $(CH_2)_nNR^{11}R^{12}$, $(CH_2)_nOR^{11}$, $(CH_2)_nO(CH_2)_kNR^{11}R^{12}$, $(CH_2)_nCOOR^{12}$, $(CH_2)_nCONR^{11}R^{12}$, $(CH_2)_nNR^{11}COR^{13}$, $(CH_2)_nNR^{11}CONR^{11}R^{12}$, $(CH_2)_nNR^{11}SO_2A$, $(CH_2)_nSO_2NR^{11}R^{12}$, $(CH_2)_nS(O)_uR^{13}$, $(CH_2)_nOC(O)R^{13}$, $(CH_2)_nCOR^{13}$, $(CH_2)_nSR^{11}$, $CH=N-OA$, $CH_2CH=N-OA$, $(CH_2)_nNHOA$, $(CH_2)_nCH=N-R^{11}$, $(CH_2)_nOC(O)NR^{11}R^{12}$, $(CH_2)_nNR^{11}COOR^{12}$, $(CH_2)_nN(R^{11})CH_2CH_2OR^{13}$, $(CH_2)_nN(R^{11})CH_2CH_2OCF_3$, $(CH_2)_nN(R^{11})C(R^{13})HCOOR^{12}$, $C(R^{13})HCOOR^{12}$, $(CH_2)_nN(R^{11})CH_2CH_2N(R^{12})CH_2COOR^{12}$, $(CH_2)_nN(R^{11})CH_2CH_2NR^{11}R^{12}$, $CH=CHCOOR^{11}$, $CH=CHCH_2NR^{11}R^{12}$, $CH=CHCH_2NR^{11}R^{12}$, $CH=CHCH_2OR^{13}$, $(CH_2)_nN(COOR^{11})COOR^{12}$, $(CH_2)_nN(CONH_2)COOR^{11}$, $(CH_2)_nN(CONH_2)CONH_2$, $(CH_2)_nN(CH_2COOR^{11})COOR^{12}$, $(CH_2)_nN(CH_2CONH_2)COOR^{11}$, $(CH_2)_nN(CH_2CONH_2)CONH_2$, $(CH_2)_nCHR^{13}COR^{11}$, $(CH_2)_nCHR^{13}COOR^{11}$, $(CH_2)_nCHR^{13}CH_2OR^{14}$, $(CH_2)_nOCN$ and $(CH_2)_nNCO$,

q is 0, 1, 2, 3, or 4,

X represents a bond or is $(CR^{11}R^{12})_h$, or $(CHR^{11})_h-Q-(CHR^{12})_i$,

Ar^2 is selected from aromatic hydrocarbons containing 6 to 14 carbon atoms and ethylenical unsaturated or aromatic heterocyclic residues containing 3 to 10 carbon atoms and one or two hetero atoms, independently selected from N, O and S,

r is 0, 1, 2, 3, 4 or 5, and

h, i are independently from each other 0, 1, 2, 3, 4, 5 or 6;

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r — is 0, 1, 2, 3, 4 or 5

and optionally isolating and/or treating the compound of formula II obtained by said reaction with an acid, to obtain the salt thereof.

29. (Cancelled):

30. (Cancelled):

31. (Currently Amended): A compound according to claim 3, wherein Ar¹ is phenyl, pyridinyl, oxazolyl, isoxazolyl, pyrazolyl or imidazolyl, preferably phenyl, pyridinyl or isoxazolyl and especially phenyl or oxazolyl,

p — is 1, 2 or 3,

R⁸ — is selected from the group consisting of alkyl comprising 1 to 4 carbon atoms, alkoxy comprising 1 to 4 carbon atoms, Hal, CH₂Hal, CH(Hal)₂, perhaloalkyl comprising 1 to 4 carbon atoms, NO₂, (CH₂)_nCN, (CH₂)_nNR¹¹R¹², (CH₂)_nCOR¹³, (CH₂)_nCOOR¹⁴, (CH₂)_nCONR¹¹R¹², (CH₂)_nSO₂NR¹¹R¹² and (CH₂)_nS(O)_uR¹³, wherein

n — is 0 or 1,

u — is 0 or 2,

q — is 0 or 1, and

X — is O, S, NR¹⁵, CHOR¹⁴, CH₂, CH₂CH₂, OCH₂, CH₂O, OCH₂CH₂, or CH₂CH₂O.

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32. (Previously Presented): A compound according to claim 31, wherein A² is phenyl or pyridinyl.
33. (Previously Presented): A compound according to claim 31, wherein X is O or S.
34. (Previously Presented): A compound according to claim 31, wherein Y is O or S.
35. (Previously Presented): A compound according to claim 31, wherein A¹ is phenyl or oxazolyl.
36. (Previously Presented): A compound according to claim 31, wherein A² is pyridinyl.
37. (Previously Presented): A compound according to claim 31, wherein X is O.
38. (Previously Presented): A compound according to claim 31, wherein Y is O.
39. (New): A compound according to claim 31, wherein A¹ is phenyl.
40. (New): A compound according to claim 3, wherein A is alkyl having 1 to 10 carbon atoms, allyl, 2-butenyl, 3-butenyl, isobutenyl, sec-butenyl, 4-pentenyl, isopentenyl, 5-hexenyl, cycloalkyl having 3 to 7 carbon atoms, alkylencycloalkyl having 5 to 10 carbon atoms, alkoxy having 1 to 10 carbon atoms, or C_uH_{2u+1}-O-(CH₂)_v wherein u is 1 to 6 and v is 1 to 6.
41. (New): A compound according to claim 4, wherein A is alkyl having 1 to 10 carbon atoms, allyl, 2-butenyl, 3-butenyl, isobutenyl, sec-butenyl, 4-pentenyl, isopentenyl, 5-hexenyl, cycloalkyl having 3 to 7 carbon atoms, alkylencycloalkyl having 5 to 10 carbon atoms, alkoxy having 1 to 10 carbon atoms, or C_uH_{2u+1}-O-(CH₂)_v wherein u is 1 to 6 and v is 1 to 6.

42. (New): A compound according to claim 31, wherein A is alkyl having 1 to 10 carbon atoms, allyl, 2-butenyl, 3-butenyl, isobutenyl, sec-butenyl, 4-pentenyl, isopentenyl, 5-hexenyl, cycloalkyl having 3 to 7 carbon atoms, alkylencycloalkyl having 5 to 10 carbon atoms, alkoxy having 1 to 10 carbon atoms, or $C_uH_{2u+1}-O-(CH_2)_v$ wherein u is 1 to 6 and v is 1 to 6.

43. (New): A compound according to claim 35, wherein A is alkyl having 1 to 10 carbon atoms, allyl, 2-butenyl, 3-butenyl, isobutenyl, sec-butenyl, 4-pentenyl, isopentenyl, 5-hexenyl, cycloalkyl having 3 to 7 carbon atoms, alkylencycloalkyl having 5 to 10 carbon atoms, alkoxy having 1 to 10 carbon atoms, or $C_uH_{2u+1}-O-(CH_2)_v$ wherein u is 1 to 6 and v is 1 to 6.

44. (New): A compound according to claim 39, wherein A is alkyl having 1 to 10 carbon atoms, allyl, 2-butenyl, 3-butenyl, isobutenyl, sec-butenyl, 4-pentenyl, isopentenyl, 5-hexenyl, cycloalkyl having 3 to 7 carbon atoms, alkylencycloalkyl having 5 to 10 carbon atoms, alkoxy having 1 to 10 carbon atoms, or $C_uH_{2u+1}-O-(CH_2)_v$ wherein u is 1 to 6 and v is 1 to 6.

45. (New): A compound according to claim 31, wherein A^2 is pyridinyl and X is O.

46. (New): A compound according to claim 35, wherein A^2 is pyridinyl and X is O.

47. (New): A compound according to claim 39, wherein A^2 is pyridinyl and X is O.

48. (New): A compound according to claim 42, wherein A^2 is pyridinyl and X is O.

49. (New): A compound according to claim 43, wherein A^2 is pyridinyl and X is O.

50. (New): A compound according to claim 44, wherein A² is pyridinyl and X is O.